```
=> screen 970 AND 2067
       SCREEN CREATED
 Ll
Uploading C:\Program Files\Stnexp\Queries\0988912.str
          STRUCTURE UPLOADED
 => que L2 AND L1
    QUE L2 AND L1
 => d
L3 HAS NO ANSWERS
Ll
                   SCR 970 AND 2067
                   STR
*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***
Structure attributes must be viewed using STN Express query preparation.
L3
                   QUE ABB=ON PLU=ON L2 AND L1
=> s 13 sss sam
SAMPLE SEARCH INITIATED 14:18:41 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED -
100.0% PROCESSED
                          29 ITERATIONS
                                                                          1 ANSWERS
SEARCH TIME: 00.00.01
FULL FILE PROJECTIONS: ONLINE **COMPLETE**
                                     **COMPLETE**
                            BATCH
PROJECTED ITERATIONS:
                                    257 TO
                                                  903
PROJECTED ANSWERS:
                                      1 TO
                                                   80
L4
                1 SEA SSS SAM L2 AND L1
=> d
     ANSWER 1 OF 1 REGISTRY COPYRIGHT 2003 ACS on STN 150048-73-0 REGISTRY 2-Propenoic acid, 2-methyl-, 1-(2-anthracenyl)ethyl ester, polymer with ethenylbenzene (9CI) (CA INDEX NAME)
L4
CN
OTHER CA INDEX NAMES:
     Benzene, ethenyl-, polymer with 1-(2-anthracenyl)ethyl 2-methyl-2-propenoate (9CI)
_{\rm CN}
MF
      (C20 H18 O2 . C8 H8)x
CI
     PMS
PCT Polyacrylic, Polystyrene
SR
     CA
     STN Files: CA, CAPLUS
LC
     CM
          1
     CRN 54720-11-5
CMF C20 H18 O2
                   СН- Ме
```

CM 2

CRN 100-42-5 CMF C8 H8

 $H_2C = CH - Ph$

- 1 REFERENCES IN FILE CA (1907 TO DATE)
- 1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

=> FIL CAPLUS HCAPLUS USPATFUL COST IN U.S. DOLLARS

SINCE FILE ENTRY 2.08

TOTAL

SESSION

FULL ESTIMATED COST

FILE 'CAPLUS' ENTERED AT 14:18:58 ON 19 NOV 2003 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2003 AMERICAN CHEMICAL SOCIETY (ACS)

FILE 'HCAPLUS' ENTERED AT 14:18:58 ON 19 NOV 2003 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2003 AMERICAN CHEMICAL SOCIETY (ACS)

FILE 'USPATFULL' ENTERED AT 14:18:58 ON 19 NOV 2003 CA INDEXING COPYRIGHT (C) 2003 AMERICAN CHEMICAL SOCIETY (ACS)

=> s 14

L5 2 L4

=> d 15 1-2 ibib hitstr

ANSWER 1 OF 2 CAPLUS COPYRIGHT 2003 ACS on STN 1993:562441 CAPLUS

ACCESSION NUMBER:

DOCUMENT NUMBER: 119:162441

TITLE: Photochemical dimerization in hydrophilicity

improvement of (in)organic articles Irie, Masahiro; Kishimoto, Soichiro INVENTOR(S):

PATENT ASSIGNEE(S): Unitika Ltd, Japan

Jpn. Kokai Tokkyo Koho, 6 pp. SOURCE:

CODEN: JKXXAF

DOCUMENT TYPE: Patent LANGUAGE: Japanese

FAMILY ACC. NUM. COUNT:

PATENT INFORMATION:

PATENT NO. KIND DATE APPLICATION NO. DATE ____ _____ JP 05024951 19930202 JP 1991-208592 19910724 PRIORITY APPLN. INFO.: JP 1991-208592 19910724

150048-73-0

RL: USES (Uses) (plates, hydrophilic treatment for with UV irradn., in presence of

RN

hydrophilic group-contg. anthracenes)
150048-73-0 CAPLUS
2-Propenoic acid, 2-methyl-, 1-(2-anthracenyl)ethyl ester, polymer with CN ethenylbenzene (9CI) (CA INDEX NAME)

CM 1

CRN 54720-11-5 CMF C20 H18 O2

CM

CRN 100-42-5 CMF C8 H8

 $H_2C = CH - Ph$

ANSWER 2 OF 2 HCAPLUS COPYRIGHT 2003 ACS on STN ACCESSION NUMBER: 1993:562441 HCAPLUS

DOCUMENT NUMBER:

TITLE:

119:162441

Photochemical dimerization in hydrophilicity

improvement of (in)organic articles Irie, Masahiro; Kishimoto, Soichiro Unitika Ltd, Japan Jpn. Kokai Tokkyo Koho, 6 pp. CODEN: JKXXAF

INVENTOR(S): PATENT ASSIGNEE(S):

SOURCE:

DOCUMENT TYPE:

Patent

LANGUAGE:

Japanese

FAMILY ACC. NUM. COUNT: PATENT INFORMATION:

PATENT NO.

KIND DATE

APPLICATION NO. DATE

JP 05024951

JP 1991-208592

19910724

PRIORITY APPLN. INFO.:

A2 19930202

JP 1991-208592

19910724

150048-73-0

RL: USES (Uses)

(plates, hydrophilic treatment for with UV irradn., in presence of

hydrophilic group-contg. anthracenes)

RN150048-73-0 HCAPLUS

2-Propenoic acid, 2-methyl-, 1-(2-anthracenyl)ethyl ester, polymer with ethenylbenzene (9CI) (CA INDEX NAME) CN

CM 1

CRN 54720-11-5 CMF C20 H18 O2

CM

CRN 100-42-5 CMF C8 H8

H2C== CH- Ph

```
1 2 3 14 16 17 18 19 21 22 56 57 63 64 66 67 68 69 70 71 73 74 75 76 77 78 80 81 82 83 84 85 87 88 89 90 91 92 94 96
ring nodes :

    4
    5
    6
    7
    8
    9
    10
    11
    12
    13
    24
    25
    26
    27
    28
    29
    30
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    32
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    41
    42
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    47
    48
    49
    50
    51
    52
    53
    54
    55

                                                                                                                        35
                                                                                                                               36
                                                                                                                                    37
chain bonds :
     1-2 1-14 1-16 2-3 3-22
16-21 17-18 17-19 24-76
39-90 40-91 41-92 44-94
                                               37-74 38-89
                                               47-88
                                                                    51 - 87
                                                         48-82
                                                                              52 - 83
ring bonds
      4-9 4-5 5-6 6-7 7-8 7-
27-28 28-29 28-30 29-33
                                          7-10
                                                 8-9
                                                          8-13
                                                                    10-11
                                                                                                                        25-26
36-37
                                                                                                                                  26-27
                                               30-31
                                                        31-32
                                                                    31 - 34
                                                                              32-33
                                                                                         32-37
                                                                                                   34 - 35
                                                                                                             35-36
                                                                                                                                  38 - 39
                                   41-42
                                                         42-44
      38-43
               39-40
                          40-41
                                               42-43
                                                                    43-47
                                                                              44-45
                                                                                         45-46
                                                                                                   45-48
                                                                                                                                  48-49
     49-50
               49-52
                          50-51
                                     50-55
                                               52-53
                                                         53-54
                                                                    54-55
exact/norm bonds :
     1-2 1-14 2-3 3-22 3-96
25-77 26-78 27-80 30-81
47-88 48-82 51-87 52-83
                                                       6-69 9-63 10-68 11-67 12-66 13-64 16-21 24-76 34-71 36-73 37-74 38-89 39-90 40-91 41-92 44-94
                                               5-70
                                               33-75
54-84
                                                                    56-57
                                                         55-85
exact bonds :
     1-16 16-17 17-18
                                   17-19
normalized bonds:
      4-9 4-5 5-6
                           6-7
                                  7-8 7-10 8-9 8-13
                                                                    10-11
                                                                              11-12
                                                                                         12-13
                                                                                                   24-25
                                                                                                             24-29
                                                                                                                        25-26
                                                                                                                                  26-27
     27-28 28-29 28-30 29-33 30-31 31-32 38-43 39-40 40-41 41-42 42-43 42-44 49-50 49-52 50-51 50-55 52-53 53-54
                                                                                                 34-35
                                                                                                             35-36
                                                                    31-34
                                                                              32-33
                                                                                         32-37
                                                                                                                        36-37
                                                                                                                                  38 - 39
                                                                    43-47
                                                                              44-45
                                                                                                                                  48-49
                                                                    54-55
```

G1:H G2:H,CH3

G3:F,H,[*1]

chain nodes :

G4:[*2],[*3],[*4]

Match level:
 1:CLASS 2:CLASS 3:CLASS 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom 12:Atom 13:Atom 14:CLASS 16:CLASS 17:CLASS 18:CLASS 19:CLASS 21:CLASS 22:CLASS 24:CLASS 25:Atom 26:Atom 27:Atom 28:Atom 29:Atom 30:Atom 31:Atom 32:Atom 33:Atom 34:Atom 35:Atom 36:Atom 37:Atom 38:Atom 39:Atom 40:Atom 41:Atom 42:Atom 43:Atom 44:Atom 45:Atom 46:Atom 47:Atom 48:Atom 49:Atom 50:Atom 51:Atom 52:Atom 53:Atom 54:Atom 55:Atom 56:CLASS 63:CLASS 64:CLASS 66:CLASS 67:CLASS 68:CLASS 69:CLASS 70:CLASS 71:CLASS 73:CLASS 74:CLASS 75:CLASS 76:CLASS 77:CLASS 78:CLASS 80:CLASS 81:CLASS 82:CLASS 83:CLASS 84:CLASS 87:CLASS 88:CLASS 89:CLASS 90:CLASS 91:CLASS 94:CLASS 96:CLASS 85:CLASS 88:CLASS 89:CLASS 88:CLASS 89:CLASS 81:CLASS 92:CLASS 94:CLASS 96:CLASS 87:CLASS 87:CLASS 88:CLASS 89:CLASS 80:CLASS 80:CLASS 92:CLASS 94:CLASS 96:CLASS 87:CLASS 87: